

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP013022.D
 Acq On : 09 Dec 2022 17:23
 Operator : CG/JU
 Sample : N5896-15
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXN3

Quant Time: Dec 09 21:56:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 08 00:35:07 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	653067	20.000	ng/u1	0.00
20) Naphthalene-d8	10.787	136	2830920	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	1838895	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	3963419	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	3076923	20.000	ng/u1	0.00
88) Perylene-d12	23.951	264	2778777	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	60721	3.975	ng/uL	0.00
4) Pyridine-d5	3.781	84	671436	15.475	ng/u1	0.00
7) Phenol-d5	7.128	99	1336799	25.131	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	781267	24.347	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	1078952	25.766	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	1112174	25.517	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	541538	26.036	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	630358	26.920	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.405	165	1192436	26.217	ng/u1	0.00
31) 4-Chloroaniline-d4	10.922	131	1376177	21.433	ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	3667683	25.952	ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	4049209	26.074	ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	432029	16.994	ng/u1	0.00
60) Fluorene-d10	15.604	176	3129746	26.176	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	488900	19.168	ng/u1	0.00
73) Anthracene-d10	17.469	188	4862642	27.195	ng/u1	0.00
81) Pyrene-d10	19.698	212	5415718	30.664	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.792	264	3805447	27.475	ng/u1	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.840	128	505366	3.413	ng/u1	99
36) 2-Methylnaphthalene	12.440	142	165910	1.633	ng/u1	97
50) Acenaphthylene	14.334	152	618517	3.705	ng/u1	99
52) Acenaphthene	14.675	153	197029	1.703	ng/u1	99
61) Fluorene	15.663	166	164976	1.251	ng/u1	98
71) Pentachlorophenol	17.016	266	239139	7.726	ng/u1	100
72) Phenanthrene	17.410	178	1986069	9.621	ng/u1	100
74) Anthracene	17.504	178	2052994	9.960	ng/u1	99
80) Fluoranthene	19.375	202	7216042	35.998	ng/u1	100
82) Pyrene	19.728	202	6452024	31.203	ng/u1	99
85) Benzo(a)anthracene	21.439	228	3257347	15.947	ng/u1	97
87) Chrysene	21.498	228	3074962	15.919	ng/u1	98
90) Benzo(b)fluoranthene	23.192	252	7090945	40.018	ng/u1	98
91) Benzo(k)fluoranthene	23.233	252	2011547	11.448	ng/u1	96
93) Benzo(a)pyrene	23.839	252	3232338	20.950	ng/u1	96
94) Indeno(1,2,3-cd)pyrene	26.551	276	2834275	14.748	ng/u1	97
95) Dibenzo(a,h)anthracene	26.563	278	706057	4.437	ng/u1#	85
96) Benzo(g,h,i)perylene	27.362	276	2213047	14.344	ng/u1	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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