

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012113.D
 Acq On : 13 Oct 2022 10:42
 Operator : CG/JU
 Sample : N5049-21
 Misc :
 ALS Vial : 84 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BE1

Quant Time: Oct 13 23:30:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	282224	20.000	ng/u1	0.00
20) Naphthalene-d8	10.651	136	1140351	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.498	164	606016	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1135518	20.000	ng/u1	0.00
79) Chrysene-d12	21.345	240	1184388	20.000	ng/u1	0.00
88) Perylene-d12	23.768	264	1278989	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	25742	3.596	ng/uL	0.00
4) Pyridine-d5	3.705	84	127771	6.712	ng/u1	0.00
7) Phenol-d5	7.016	99	165254	7.281	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	418694	32.498	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	476675	26.457	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	293310	16.824	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	289636	35.963	ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	197158	24.143	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	451865	28.551	ng/u1	0.00
31) 4-Chloroaniline-d4	10.799	131	524378	22.484	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	1359794	35.430	ng/u1	0.00
49) Acenaphthylene-d8	14.192	160	1632417	34.396	ng/u1	0.00
54) 4-Nitrophenol-d4	14.728	143	11255	1.494	ng/u1	0.02
60) Fluorene-d10	15.492	176	1209336	34.827	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	46831	7.811	ng/u1	0.00
73) Anthracene-d10	17.357	188	1848499	37.488	ng/u1	0.00
81) Pyrene-d10	19.592	212	2150935	35.449	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.610	264	2363975	38.022	ng/u1	0.00

Target Compounds Qvalue

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

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