

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012992.D
 Acq On : 08 Dec 2022 19:23
 Operator : CG/JU
 Sample : N5894-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 09 01:58:13 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	551454	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	2400406	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164	1592311	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.369	188	3638174	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	3444249	20.000	ng/ul	0.00
88) Perylene-d12	23.945	264	3646773	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	61692	4.783	ng/uL	0.00
4) Pyridine-d5	3.787	84	193753	5.288	ng/ul	0.00
7) Phenol-d5	7.128	99	305843	6.809	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	834437	30.796	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	858931	24.292	ng/ul	0.00
15) 4-Methylphenol-d8	8.681	113	595178	16.172	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	540639	30.655	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	574080	28.913	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.405	165	1085844	28.155	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	681270	12.513	ng/ul	0.00
46) Dimethylphthalate-d6	14.022	166	4100747	33.509	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160	4353572	32.376	ng/ul	0.00
54) 4-Nitrophenol-d4	14.804	143	103865	4.718	ng/ul	0.00
60) Fluorene-d10	15.604	176	3567771	34.461	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	583455	24.921	ng/ul	0.00
73) Anthracene-d10	17.469	188	5866841	35.744	ng/ul	0.00
81) Pyrene-d10	19.698	212	7050334	35.662	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.786	264	6513678	35.835	ng/ul	0.00

Target Compounds Qvalue

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

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