

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
 Data File : BP019357.D
 Acq On : 13 Jan 2024 07:09
 Operator : MA/JU
 Sample : P1091-20
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B03

Quant Time: Jan 14 20:26:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 09 03:05:24 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	135085	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.569	136	527824	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.422	164	373363	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.181	188	949295	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	965594	20.000	ng/u1	0.00
88) Perylene-d12	23.639	264	971016	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	14868	4.162	ng/uL	0.00
4) Pyridine-d5	3.640	84	100238	9.703	ng/u1	0.00
7) Phenol-d5	6.964	99	80294	6.219	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.117	67	285616	36.984	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.311	132	240192	27.082	ng/u1	0.00
15) 4-Methylphenol-d8	8.505	113	159070	15.530	ng/u1	-0.01
21) Nitrobenzene-d5	8.946	128	170934	38.494	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	191298	37.076	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.205	165	343745	32.977	ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	423682	32.177	ng/u1	0.00
46) Dimethylphthalate-d6	13.840	166	1288957	39.076	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	1350273	38.188	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	27619	5.453	ng/u1	0.00
60) Fluorene-d10	15.416	176	1145098	42.402	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	242060	34.538	ng/u1	0.00
73) Anthracene-d10	17.281	188	2042916	43.149	ng/u1	0.00
81) Pyrene-d10	19.528	212	2657651	42.210	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.492	264	2359026	44.680	ng/u1	0.00

Target Compounds Qvalue

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

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