

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
 Data File : BP018566.D
 Acq On : 04 Dec 2023 16:36
 Operator : MA/JU
 Sample : 05441-04
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0QF9

Quant Time: Dec 04 17:09:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 01 21:56:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	308929	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.651	136	1355754	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.486	164	896846	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1841219	20.000	ng/u1	-0.01
79) Chrysene-d12	21.316	240	1169392	20.000	ng/u1	-0.01
88) Perylene-d12	23.692	264	1130608	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	31385	3.462	ng/uL	0.00
4) Pyridine-d5	3.687	84	196821	8.105	ng/u1	0.00
7) Phenol-d5	7.022	99	141527	4.565	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.193	67	541864	29.490	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.387	132	458277	18.944	ng/u1	-0.01
15) 4-Methylphenol-d8	8.563	113	272550	10.863	ng/u1	-0.02
21) Nitrobenzene-d5	9.016	128	319187	26.452	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.740	143	300656	24.033	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.269	165	576275	22.528	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.793	131	586876	17.555	ng/u1	-0.01
46) Dimethylphthalate-d6	13.904	166	2363727	29.530	ng/u1	0.00
49) Acenaphthylene-d8	14.181	160	2486821	28.215	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.681	143	38377	2.921	ng/u1	-0.01
60) Fluorene-d10	15.475	176	2025058	30.140	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.592	200	249732	22.621	ng/u1	-0.01
73) Anthracene-d10	17.333	188	3053452	31.545	ng/u1	0.00
81) Pyrene-d10	19.569	212	3223313	35.410	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.539	264	2223665	33.732	ng/u1	-0.01

Target Compounds Qvalue

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

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