

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018384.D
 Acq On : 26 Nov 2023 00:22
 Operator : MA/JU
 Sample : 05261-11
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 27 00:05:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 24 22:02:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	584192	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	2354480	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	1451670	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	2775316	20.000	ng/ul	0.00
79) Chrysene-d12	21.351	240	2041130	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	2517136	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	49764	3.191	ng/uL	0.00
4) Pyridine-d5	3.687	84	634183	14.795	ng/ul	0.00
7) Phenol-d5	7.028	99	1135053	21.260	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	605715	18.973	ng/ul	0.00
11) 2-Chlorophenol-d4	7.393	132	925381	20.644	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	1068047	24.789	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	520113	27.197	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	571285	32.346	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	1212893	28.465	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	1564123	27.217	ng/ul	0.00
46) Dimethylphthalate-d6	13.916	166	4380553	33.913	ng/ul	0.00
49) Acenaphthylene-d8	14.192	160	4726814	32.115	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	479498	26.890	ng/ul	0.00
60) Fluorene-d10	15.492	176	3574722	33.087	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	400275	33.880	ng/ul	0.00
73) Anthracene-d10	17.345	188	5061034	34.118	ng/ul	0.00
81) Pyrene-d10	19.586	212	5221951	32.145	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	5324526	36.486	ng/ul	0.02

Target Compounds Qvalue

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

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