

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018406.D
 Acq On : 27 Nov 2023 18:21
 Operator : MA/JU
 Sample : 05341-15
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 27 21:58:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 27 21:50:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	498762	20.000	ng/ul	0.00
20) Naphthalene-d8	10.658	136	1992662	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.493	164	1123205	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	1959854	20.000	ng/ul	0.00
79) Chrysene-d12	21.375	240	1553804	20.000	ng/ul	0.00
88) Perylene-d12	23.886	264	1871683	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	48734	3.660	ng/uL	0.00
4) Pyridine-d5	3.687	84	211512	5.780	ng/ul	0.00
7) Phenol-d5	7.022	99	253625	5.564	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	684959	25.130	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	747520	19.532	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	459993	12.505	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	466928	28.849	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	433061	28.972	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	838340	23.247	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	944488	19.419	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	2818628	28.202	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	3148067	27.644	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	42165	3.056	ng/ul	0.00
60) Fluorene-d10	15.487	176	2356820	28.193	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	206544	24.757	ng/ul	0.00
73) Anthracene-d10	17.345	188	3434510	32.787	ng/ul	0.00
81) Pyrene-d10	19.592	212	3446662	27.871	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.722	264	3690889	34.014	ng/ul	0.02

Target Compounds Qvalue

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

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