

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018405.D
 Acq On : 27 Nov 2023 17:47
 Operator : MA/JU
 Sample : 05341-14
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 27 21:57:52 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	7.863	152	448184	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	1704364	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	912787	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	1619301	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	1537918	20.000	ng/ul	0.00
88) Perylene-d12	23.880	264	1814184	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	42050	3.514	ng/uL	0.00
4) Pyridine-d5	3.687	84	246208	7.487	ng/ul	0.00
7) Phenol-d5	7.022	99	190740	4.657	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	591878	24.165	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	628042	18.262	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	360355	10.902	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	386584	27.925	ng/ul	0.00
24) 2-Nitrophenol-d4	9.745	143	344400	26.938	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	658276	21.342	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	853412	20.514	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	2079552	25.604	ng/ul	0.00
49) Acenaphthylene-d8	14.186	160	2379641	25.713	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	28541	2.545	ng/ul	0.00
60) Fluorene-d10	15.486	176	1722283	25.352	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	134923	19.573	ng/ul	0.00
73) Anthracene-d10	17.345	188	2464408	28.474	ng/ul	0.00
81) Pyrene-d10	19.592	212	2781821	22.727	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.715	264	3019759	28.711	ng/ul	0.02

Target Compounds Qvalue

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

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