

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120522\
 Data File : BP012955.D
 Acq On : 05 Dec 2022 16:39
 Operator : CG/JU
 Sample : N5738-18DL 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBNJ7DL

Quant Time: Dec 05 21:52:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 01 23:46:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	568894	20.000	ng/u1	0.00
20) Naphthalene-d8	10.822	136	2560515	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.640	164	1776235	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	3955021	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	3077840	20.000	ng/u1	0.00
88) Perylene-d12	23.992	264	2775794	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	4332	0.339	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.158	99	69926	1.331	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.322	67	53513	1.799	ng/u1	0.00
11) 2-Chlorophenol-d4	7.528	132	62148	1.614	ng/u1	0.00
15) 4-Methylphenol-d8	8.705	113	45021	1.044	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	29355	1.612	ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	29333	1.408	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	62156	1.490	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	24777	0.420	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	257430	2.009	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	271418	1.919	ng/u1	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/u1	
60) Fluorene-d10	15.634	176	245478	2.180	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/u1	
73) Anthracene-d10	17.498	188	392069	2.243	ng/u1	0.00
81) Pyrene-d10	19.722	212	437285	2.890	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.821	264	262618	1.936	ng/u1	0.00
Target Compounds						
					Qvalue	
71) Pentachlorophenol	17.039	266	1371336	43.296	ng/u1	99

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

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