

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012372.D
 Acq On : 28 Oct 2022 11:48
 Operator : CG/JU
 Sample : N5232-11
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA65

Quant Time: Oct 28 23:20:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 28 06:05:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	250403	20.000	ng/u1	0.00
20) Naphthalene-d8	10.604	136	1161417	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.457	164	779435	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.216	188	1765201	20.000	ng/u1	0.00
79) Chrysene-d12	21.310	240	2000740	20.000	ng/u1	0.00
88) Perylene-d12	23.704	264	1859081	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	21351	3.433	ng/uL	0.00
4) Pyridine-d5	3.699	84	16393	0.918	ng/u1	0.03
7) Phenol-d5	6.981	99	145582	6.482	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	407924	30.425	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	412013	25.121	ng/u1	0.00
15) 4-Methylphenol-d8	8.522	113	291164	16.391	ng/u1	0.00
21) Nitrobenzene-d5	8.975	128	281999	37.442	ng/u1	0.00
24) 2-Nitrophenol-d4	9.693	143	296072	37.874	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	534235	31.534	ng/u1	0.00
31) 4-Chloroaniline-d4	10.763	131	51490	2.041	ng/u1	0.01
46) Dimethylphthalate-d6	13.869	166	2074901	38.887	ng/u1	0.00
49) Acenaphthylene-d8	14.151	160	2270864	37.243	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	62793	6.218	ng/u1	0.01
60) Fluorene-d10	15.451	176	1815415	39.740	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	317011	34.494	ng/u1	0.00
73) Anthracene-d10	17.316	188	3160683	41.556	ng/u1	0.00
81) Pyrene-d10	19.557	212	3945690	39.280	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.551	264	3840771	42.282	ng/u1	0.00

Target Compounds	Qvalue

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

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