

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
 Data File : BP014402.D
 Acq On : 30 Mar 2023 15:19
 Operator : CG/JU
 Sample : 02059-01DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0QG0DL

Quant Time: Mar 30 22:40:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 30 12:32:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	159791	20.000	ng/u1	0.00
20) Naphthalene-d8	10.840	136	713859	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.669	164	492856	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.433	188	1093899	20.000	ng/u1	0.00
79) Chrysene-d12	21.521	240	862588	20.000	ng/u1	0.00
88) Perylene-d12	24.045	264	782269	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	4286	1.040	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.193	99	16441	1.116	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.340	67	57550	6.087	ng/u1	0.00
11) 2-Chlorophenol-d4	7.552	132	47954	4.502	ng/u1	0.00
15) 4-Methylphenol-d8	8.746	113	32162	2.706	ng/u1	0.01
21) Nitrobenzene-d5	9.199	128	31024	5.383	ng/u1	0.00
24) 2-Nitrophenol-d4	9.928	143	32531	4.930	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.475	165	56185	4.675	ng/u1	0.01
31) 4-Chloroaniline-d4	10.999	131	51572	3.258	ng/u1	0.01
46) Dimethylphthalate-d6	14.075	166	289669	7.249	ng/u1	0.00
49) Acenaphthylene-d8	14.363	160	287302	6.479	ng/u1	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/u1	
60) Fluorene-d10	15.669	176	242336	7.272	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	36242	4.503	ng/u1	0.00
73) Anthracene-d10	17.533	188	385900	7.289	ng/u1	0.00
81) Pyrene-d10	19.763	212	434913	7.506	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.880	264	297384	7.155	ng/u1	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.893	128	1182258	29.733	ng/u1	98
36) 2-Methylnaphthalene	12.498	142	148317	5.504	ng/u1	99
37) 1-Methylnaphthalene	12.716	142	118007	4.416	ng/u1	98
43) 1,1'-Biphenyl	13.504	154	46798	1.175	ng/u1	97
58) 2,3,4,6-Tetrachlorophenol	15.304	232	78782	7.572	ng/u1	98
71) Pentachlorophenol	17.086	266	1540609	170.082	ng/u1	99
72) Phenanthrene	17.475	178	92841	1.526	ng/u1	98

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

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