

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014229.D
 Acq On : 22 Mar 2023 22:11
 Operator : CG/JU
 Sample : 01977-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYG92

Quant Time: Mar 23 02:36:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 23 02:31:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	162327	20.000	ng/u1	0.00
20) Naphthalene-d8	10.863	136	735949	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.687	164	541244	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	1271977	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.527	240	1303499	20.000	ng/u1	# 0.00
88) Perylene-d12	24.045	264	1341943	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	30674	7.140	ng/uL	0.00
4) Pyridine-d5	3.840	84	417387	35.088	ng/u1	0.00
7) Phenol-d5	7.199	99	579657	40.465	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	381053	45.611	ng/u1	0.00
11) 2-Chlorophenol-d4	7.569	132	438087	40.140	ng/u1	0.00
15) 4-Methylphenol-d8	8.752	113	479638	40.991	ng/u1	0.00
21) Nitrobenzene-d5	9.216	128	237177	40.390	ng/u1	0.00
24) 2-Nitrophenol-d4	9.940	143	285555	40.624	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.487	165	482675	36.494	ng/u1	0.00
31) 4-Chloroaniline-d4	11.004	131	658969	37.649	ng/u1	0.00
46) Dimethylphthalate-d6	14.092	166	1819893	39.620	ng/u1	0.00
49) Acenaphthylene-d8	14.381	160	1943532	39.941	ng/u1	0.00
54) 4-Nitrophenol-d4	14.892	143	293269	36.230	ng/u1	0.00
60) Fluorene-d10	15.681	176	1533153	39.383	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.804	200	321557	33.448	ng/u1	0.00
73) Anthracene-d10	17.545	188	2501678	40.514	ng/u1	0.00
81) Pyrene-d10	19.769	212	3046579	42.377	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.886	264	2923358	40.954	ng/u1	0.00

Target Compounds Qvalue

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

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