

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013304.D
 Acq On : 24 Dec 2022 17:31
 Operator : CG/JU
 Sample : N6016-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYD5

Quant Time: Dec 25 00:22:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 24 00:19:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	560272	20.000	ng/u1	0.00
20) Naphthalene-d8	10.751	136	2090073	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.586	164	1131917	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	2500558	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	2712546	20.000	ng/u1	0.00
88) Perylene-d12	23.915	264	3148825	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	88138	6.726	ng/uL	0.00
4) Pyridine-d5	3.758	84	1060812	28.498	ng/u1	0.00
7) Phenol-d5	7.105	99	1525040	33.418	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	1007976	36.615	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	1291297	35.945	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	966683	25.852	ng/u1	0.00
21) Nitrobenzene-d5	9.110	128	620604	40.414	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	698957	40.430	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	1176624	35.039	ng/u1	0.00
31) 4-Chloroaniline-d4	10.893	131	1128083	23.796	ng/u1	0.00
46) Dimethylphthalate-d6	13.992	166	3084893	35.461	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	3641362	38.093	ng/u1	0.00
54) 4-Nitrophenol-d4	14.787	143	486837	31.111	ng/u1	0.00
60) Fluorene-d10	15.581	176	2726540	37.047	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	368214	22.882	ng/u1	0.00
73) Anthracene-d10	17.445	188	4021524	35.648	ng/u1	0.00
81) Pyrene-d10	19.680	212	5638339	36.213	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.757	264	5870983	37.407	ng/u1	0.00

Target Compounds Qvalue

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

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