

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013196.D
 Acq On : 21 Dec 2022 08:24
 Operator : CG/JU
 Sample : N5984-13
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COLE4

Quant Time: Dec 21 21:50:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 19 21:43:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	504345	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	2076689	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	1223071	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	2288995	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	2070146	20.000	ng/ul	0.00
88) Perylene-d12	23.916	264	2465344	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	60977	5.169	ng/uL	0.00
4) Pyridine-d5	3.764	84	572874	17.097	ng/ul	0.00
7) Phenol-d5	7.111	99	1186935	28.893	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	762534	30.771	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	964929	29.838	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	960781	28.544	ng/ul	-0.01
21) Nitrobenzene-d5	9.122	128	480136	31.468	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	547799	31.890	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	979097	29.345	ng/ul	0.00
31) 4-Chloroaniline-d4	10.904	131	1172543	24.893	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	2878534	30.623	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	3340615	32.343	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	328760	19.444	ng/ul	0.00
60) Fluorene-d10	15.587	176	2470693	31.069	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	286132	19.425	ng/ul	0.00
73) Anthracene-d10	17.451	188	3415411	33.074	ng/ul	0.00
81) Pyrene-d10	19.680	212	3792996	31.921	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.757	264	4070771	33.127	ng/ul	0.00

Target Compounds Qvalue

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

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