

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018308.D
 Acq On : 23 Nov 2023 10:28
 Operator : MA/JU
 Sample : 05390-03
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0TJ5

Quant Time: Nov 23 23:50:42 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 22 14:09:55 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	373367	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	1580894	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	1035193	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	2104748	20.000	ng/u1	0.00
79) Chrysene-d12	21.357	240	1267961	20.000	ng/u1	0.00
88) Perylene-d12	23.780	264	1274555	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	51591	5.175	ng/uL	0.00
4) Pyridine-d5	3.687	84	405112	14.788	ng/u1	0.00
7) Phenol-d5	7.028	99	331580	9.717	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	733804	35.963	ng/u1	0.00
11) 2-Chlorophenol-d4	7.399	132	747196	26.081	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	525992	19.101	ng/u1	-0.01
21) Nitrobenzene-d5	9.034	128	497102	38.713	ng/u1	0.00
24) 2-Nitrophenol-d4	9.758	143	454049	38.288	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	890972	31.142	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.810	131	1397161	36.208	ng/u1	0.00
46) Dimethylphthalate-d6	13.922	166	3749900	40.710	ng/u1	-0.01
49) Acenaphthylene-d8	14.198	160	4086143	38.932	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	93559	7.358	ng/u1	-0.01
60) Fluorene-d10	15.498	176	3120494	40.503	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	270075	30.143	ng/u1	0.00
73) Anthracene-d10	17.357	188	4614209	41.016	ng/u1	0.00
81) Pyrene-d10	19.592	212	4529279	44.882	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.622	264	3137584	42.461	ng/u1	0.00

Target Compounds Qvalue

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

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