

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017490.D
 Acq On : 02 Oct 2023 22:32
 Operator : MA/JU
 Sample : 04588-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBKH0

Quant Time: Oct 02 23:17:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 04:02:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	135989	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	541726	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.640	164	323351	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	713932	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	685415	20.000	ng/ul	0.00
88) Perylene-d12	24.121	264	761885	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	13391	4.045	ng/uL	0.00
4) Pyridine-d5	3.734	84	84561	9.136	ng/ul	0.00
7) Phenol-d5	7.099	99	69743	6.244	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	263580	39.105	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	247060	27.890	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	141926	16.345	ng/ul	0.00
21) Nitrobenzene-d5	9.152	128	170594	43.559	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	173738	40.856	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	293252	36.501	ng/ul	0.00
31) 4-Chloroaniline-d4	10.952	131	364969	31.360	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	1060054	45.763	ng/ul	0.00
49) Acenaphthylene-d8	14.340	160	1169864	43.912	ng/ul	0.00
54) 4-Nitrophenol-d4	14.887	143	13344	3.307	ng/ul	0.01
60) Fluorene-d10	15.640	176	917361	46.002	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	135086	33.571	ng/ul	0.00
73) Anthracene-d10	17.528	188	1513776	47.808	ng/ul	0.00
81) Pyrene-d10	19.775	212	1858344	50.137	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	1774842	48.384	ng/ul	0.00

Target Compounds Qvalue

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

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