

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017030.D
 Acq On : 09 Sep 2023 02:00
 Operator : MA/JU
 Sample : 04166-06
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHH1

Quant Time: Sep 09 04:21:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 09 01:41:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	426031	20.000	ng/u1	0.00
20) Naphthalene-d8	10.798	136	1865597	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	1122183	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	2438510	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	2016052	20.000	ng/u1	0.00
88) Perylene-d12	24.027	264	2017087	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	47539	4.221	ng/uL	0.00
4) Pyridine-d5	3.764	84	256184	7.960	ng/u1	0.00
7) Phenol-d5	7.122	99	338736	8.703	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	1035487	42.532	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	915406	32.496	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	627388	21.158	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	593859	41.031	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	626997	39.963	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	1045661	36.100	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	542232	12.975	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	3662216	43.842	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	4039252	41.900	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	110771	6.789	ng/u1	0.00
60) Fluorene-d10	15.628	176	3091013	43.219	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	516637	35.206	ng/u1	0.00
73) Anthracene-d10	17.498	188	4868009	43.470	ng/u1	0.00
81) Pyrene-d10	19.733	212	5524139	44.986	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.862	264	4660083	45.539	ng/u1	0.00

Target Compounds	Qvalue

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

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