

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017018.D
 Acq On : 08 Sep 2023 18:31
 Operator : MA/JU
 Sample : 04166-18
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJY0

Quant Time: Sep 09 00:45:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	563163	20.000	ng/u1	0.00
20) Naphthalene-d8	10.805	136	2439639	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	1461482	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	3035609	20.000	ng/u1	-0.01
79) Chrysene-d12	21.498	240	2215262	20.000	ng/u1	-0.01
88) Perylene-d12	24.033	264	2050517	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	71919	4.831	ng/uL	0.00
4) Pyridine-d5	3.764	84	461238	10.842	ng/u1	0.00
7) Phenol-d5	7.122	99	467062	9.078	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	1315458	40.875	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	1198927	32.197	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	843273	21.514	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	767000	40.524	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	821364	40.033	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.410	165	1339052	35.351	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	1818442	33.275	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	4695871	43.165	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	5166550	41.152	ng/u1	0.00
54) 4-Nitrophenol-d4	14.846	143	168084	7.910	ng/u1	-0.01
60) Fluorene-d10	15.628	176	3943197	42.335	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	685173	37.506	ng/u1	-0.01
73) Anthracene-d10	17.504	188	6145602	44.084	ng/u1	0.00
81) Pyrene-d10	19.739	212	6547765	48.527	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.869	264	4892882	47.035	ng/u1	-0.01

Target Compounds	Qvalue

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

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