

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017016.D
 Acq On : 08 Sep 2023 17:22
 Operator : MA/JU
 Sample : 04166-16
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJP0

Quant Time: Sep 09 01:13:03 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	461459	20.000	ng/u1	0.00
20) Naphthalene-d8	10.799	136	2026000	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.634	164	1236634	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	2590683	20.000	ng/u1	-0.01
79) Chrysene-d12	21.498	240	1960547	20.000	ng/u1	-0.01
88) Perylene-d12	24.033	264	1853844	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	64237	5.265	ng/uL	0.00
4) Pyridine-d5	3.764	84	312378	8.961	ng/u1	0.00
7) Phenol-d5	7.122	99	383736	9.102	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.317	67	1030656	39.084	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	938808	30.768	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	671920	20.920	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	601426	38.264	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	634838	37.259	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.410	165	1041966	33.124	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	1445467	31.850	ng/u1	0.00
46) Dimethylphthalate-d6	14.046	166	3687428	40.058	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	4037774	38.008	ng/u1	0.00
54) 4-Nitrophenol-d4	14.846	143	142533	7.927	ng/u1	-0.01
60) Fluorene-d10	15.628	176	3178133	40.325	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	548419	35.176	ng/u1	0.00
73) Anthracene-d10	17.504	188	5184167	43.574	ng/u1	0.00
81) Pyrene-d10	19.739	212	5632382	47.166	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.869	264	4295424	45.672	ng/u1	-0.01

Target Compounds Qvalue

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

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