

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017010.D
 Acq On : 08 Sep 2023 13:55
 Operator : MA/JU
 Sample : 04123-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH8W9

Quant Time: Sep 09 01:04:15 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	547709	20.000	ng/u1	0.00
20) Naphthalene-d8	10.805	136	2387902	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	1453929	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	3089936	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	2606843	20.000	ng/u1	0.00
88) Perylene-d12	24.045	264	2630272	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	81054	5.598	ng/uL	0.00
4) Pyridine-d5	3.764	84	511769	12.369	ng/u1	0.00
7) Phenol-d5	7.122	99	444860	8.890	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	1101804	35.202	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	1035892	28.604	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	760625	19.953	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	642246	34.668	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	689370	34.328	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.410	165	1154946	31.152	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	1340930	25.069	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	3837809	35.461	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	4379393	35.063	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	154178	7.293	ng/u1	-0.01
60) Fluorene-d10	15.634	176	3305756	35.675	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	563468	30.302	ng/u1	0.00
73) Anthracene-d10	17.504	188	5196129	36.618	ng/u1	0.00
81) Pyrene-d10	19.745	212	5839475	36.777	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.880	264	5006318	37.518	ng/u1	0.00

Target Compounds Qvalue

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

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