

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017011.D
 Acq On : 08 Sep 2023 14:29
 Operator : MA/JU
 Sample : 04123-13
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH9Y1

Quant Time: Sep 09 00:44:05 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	435473	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1896251	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	1167893	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	2521688	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	2050805	20.000	ng/u1	0.00
88) Perylene-d12	24.045	264	2046638	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	66597	5.785	ng/uL	0.00
4) Pyridine-d5	3.764	84	365280	11.104	ng/u1	0.00
7) Phenol-d5	7.122	99	303348	7.625	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	928471	37.310	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	808295	28.072	ng/u1	-0.01
15) 4-Methylphenol-d8	8.681	113	562574	18.561	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	547183	37.195	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	560383	35.140	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.410	165	949876	32.263	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	1140965	26.861	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	3369335	38.757	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	3790381	37.780	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	97343	5.732	ng/u1	-0.01
60) Fluorene-d10	15.634	176	2882208	38.722	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	436118	28.739	ng/u1	0.00
73) Anthracene-d10	17.504	188	4495159	38.816	ng/u1	0.00
81) Pyrene-d10	19.745	212	5067438	40.567	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.874	264	4244332	40.878	ng/u1	0.00

Target Compounds Qvalue

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

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