

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017035.D
 Acq On : 09 Sep 2023 04:52
 Operator : MA/JU
 Sample : 04166-14
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJX8

Quant Time: Sep 09 05:35:43 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	7.969	152	327739	20.000	ng/u1	0.00
20) Naphthalene-d8	10.799	136	1453660	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	893819	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	1935888	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	1524577	20.000	ng/u1	0.00
88) Perylene-d12	24.027	264	1495404	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	41680	4.810	ng/uL	0.00
4) Pyridine-d5	3.764	84	332994	13.450	ng/u1	0.00
7) Phenol-d5	7.122	99	301929	10.084	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	806916	43.084	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	743245	34.298	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	535385	23.471	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	463723	41.119	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	498168	40.750	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	828701	36.717	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	1118369	34.345	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	2922556	43.926	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	3211987	41.832	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	100747	7.752	ng/u1	0.00
60) Fluorene-d10	15.628	176	2475861	43.463	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	415110	35.632	ng/u1	0.00
73) Anthracene-d10	17.498	188	4012051	45.128	ng/u1	0.00
81) Pyrene-d10	19.733	212	4458378	48.011	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.863	264	3523244	46.441	ng/u1	0.00

Target Compounds Qvalue

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

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