

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017029.D
 Acq On : 09 Sep 2023 01:26
 Operator : MA/JU
 Sample : 04153-08
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Sep 09 04:21:26 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	7.975	152	375543	20.000	ng/u1	0.00
20) Naphthalene-d8	10.799	136	1643625	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	985174	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	1953966	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	1433123	20.000	ng/u1	0.00
88) Perylene-d12	24.027	264	1348936	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	50621	5.099	ng/uL	0.00
4) Pyridine-d5	3.764	84	382565	13.486	ng/u1	0.00
7) Phenol-d5	7.122	99	315831	9.205	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.311	67	844409	39.347	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	763149	30.733	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	554828	21.227	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	477957	37.483	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	493398	35.695	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	833144	32.648	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	1239132	33.656	ng/u1	0.00
46) Dimethylphthalate-d6	14.046	166	3000571	40.917	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	3191387	37.709	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	91820	6.410	ng/u1	0.00
60) Fluorene-d10	15.628	176	2459082	39.165	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	357164	30.374	ng/u1	0.00
73) Anthracene-d10	17.498	188	3761186	41.915	ng/u1	0.00
81) Pyrene-d10	19.733	212	4058384	46.492	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.863	264	2954266	43.170	ng/u1	0.00

Target Compounds Qvalue

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

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