

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017076.D
 Acq On : 10 Sep 2023 05:32
 Operator : MA/JU
 Sample : 04227-13
 Misc :
 ALS Vial : 88 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJW6

Quant Time: Sep 10 06:03:55 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	567882	20.000	ng/u1	0.00
20) Naphthalene-d8	10.793	136	2473022	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.628	164	1465002	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.392	188	3097356	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	2251331	20.000	ng/u1	0.00
88) Perylene-d12	24.021	264	2163167	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	74332	4.951	ng/uL	0.00
4) Pyridine-d5	3.764	84	639252	14.902	ng/u1	0.00
7) Phenol-d5	7.122	99	434112	8.367	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	1351628	41.650	ng/u1	0.00
11) 2-Chlorophenol-d4	7.487	132	1229003	32.731	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	825905	20.896	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	792598	41.311	ng/u1	0.00
24) 2-Nitrophenol-d4	9.881	143	844313	40.596	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	1424657	37.104	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	1794696	32.397	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	4808221	44.091	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	5504290	43.736	ng/u1	0.00
54) 4-Nitrophenol-d4	14.851	143	147764	6.937	ng/u1	0.00
60) Fluorene-d10	15.622	176	4101823	43.932	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	660986	35.461	ng/u1	0.00
73) Anthracene-d10	17.498	188	6260589	44.013	ng/u1	0.00
81) Pyrene-d10	19.733	212	6688544	48.776	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.857	264	4966670	45.258	ng/u1	0.00

Target Compounds Qvalue

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

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