

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017326.D
 Acq On : 25 Sep 2023 18:09
 Operator : MA/JU
 Sample : PB155714BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK714

Quant Time: Sep 25 18:53:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	177404	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.752	136	756703	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.593	164	466089	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.363	188	1010712	20.000	ng/u1	0.00
79) Chrysene-d12	21.469	240	914838	20.000	ng/u1	0.00
88) Perylene-d12	23.986	264	980302	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	28902	6.692	ng/uL	-0.01
4) Pyridine-d5	3.728	84	416587	34.501	ng/u1	-0.01
7) Phenol-d5	7.081	99	481188	33.026	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.263	67	301851	34.329	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.446	132	393333	34.037	ng/u1	-0.02
15) 4-Methylphenol-d8	8.640	113	386973	34.162	ng/u1	-0.02
21) Nitrobenzene-d5	9.122	128	187789	34.328	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.840	143	208911	35.170	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.363	165	369864	32.958	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.916	131	552649	33.996	ng/u1	-0.01
46) Dimethylphthalate-d6	14.004	166	1212068	36.301	ng/u1	-0.01
49) Acenaphthylene-d8	14.293	160	1299559	33.841	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	157253	27.033	ng/u1	0.00
60) Fluorene-d10	15.587	176	1014559	35.295	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.722	200	123466	21.674	ng/u1	0.00
73) Anthracene-d10	17.463	188	1620620	36.154	ng/u1	0.00
81) Pyrene-d10	19.710	212	1877268	37.946	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.821	264	1783159	37.780	ng/u1	-0.01

Target Compounds	Qvalue

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

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