

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017311.D
 Acq On : 25 Sep 2023 09:28
 Operator : MA/JU
 Sample : PB155714BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK714

Quant Time: Sep 25 11:09:57 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	7.928	152	164775	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.751	136	674889	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.592	164	394028	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.363	188	890186	20.000	ng/u1	0.00
79) Chrysene-d12	21.468	240	848836	20.000	ng/u1	0.00
88) Perylene-d12	23.986	264	914480	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	27503	6.856	ng/uL	0.00
4) Pyridine-d5	3.734	84	389136	34.698	ng/u1	0.00
7) Phenol-d5	7.081	99	449437	33.211	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.269	67	272811	33.404	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.451	132	366124	34.111	ng/u1	-0.01
15) 4-Methylphenol-d8	8.640	113	346396	32.923	ng/u1	-0.02
21) Nitrobenzene-d5	9.128	128	164534	33.723	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.839	143	177619	33.527	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.369	165	329204	32.891	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.916	131	485766	33.504	ng/u1	-0.01
46) Dimethylphthalate-d6	14.004	166	975723	34.567	ng/u1	-0.01
49) Acenaphthylene-d8	14.292	160	1079978	33.266	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	129645	26.363	ng/u1	0.00
60) Fluorene-d10	15.586	176	860002	35.390	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.721	200	99819	19.895	ng/u1	0.00
73) Anthracene-d10	17.463	188	1427932	36.168	ng/u1	0.00
81) Pyrene-d10	19.704	212	1723824	37.554	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.821	264	1649587	37.466	ng/u1	-0.01

Target Compounds Qvalue

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

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