

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032524\
 Data File : BP019983.D
 Acq On : 25 Mar 2024 15:59
 Operator : MA/JU
 Sample : PB159444BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK444

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 03/26/2024
 Supervised By :mohammad ahmed 03/27/2024

Quant Time: Mar 26 00:10:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	136308m	20.000	ng/u1	-0.05
20) Naphthalene-d8	10.328	136	592085	20.000	ng/u1	-0.04
38) Acenaphthene-d10	14.228	164	435652	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.069	188	939817	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	749528	20.000	ng/u1	0.01
88) Perylene-d12	24.856	264	674516	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	22231m	7.146	ng/uL	-0.02
4) Pyridine-d5	3.610	84	326958	34.806	ng/u1	-0.03
7) Phenol-d5	6.757	99	391780	34.079	ng/u1	-0.04
9) Bis-(2-Chloroethyl)eth...	6.934	67	279843	37.363	ng/u1	-0.04
11) 2-Chlorophenol-d4	7.110	132	285456	34.576	ng/u1	-0.04
15) 4-Methylphenol-d8	8.269	113	323950	36.422	ng/u1	-0.04
21) Nitrobenzene-d5	8.722	128	156370	34.456	ng/u1	-0.04
24) 2-Nitrophenol-d4	9.434	143	182113	35.866	ng/u1	-0.04
28) 2,4-Dichlorophenol-d3	9.963	165	302115	32.579	ng/u1	-0.04
31) 4-Chloroaniline-d4	10.475	131	443330	36.673	ng/u1	-0.04
46) Dimethylphthalate-d6	13.633	166	1168978	38.271	ng/u1	-0.02
49) Acenaphthylene-d8	13.916	160	1225566	34.348	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.469	143	164565m	34.271	ng/u1	-0.01
60) Fluorene-d10	15.239	176	949887	36.507	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.398	200	193442	34.386	ng/u1	-0.02
73) Anthracene-d10	17.169	188	1538685	36.986	ng/u1	0.00
81) Pyrene-d10	19.568	212	1733380	39.417	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.621	264	1256528	38.776	ng/u1	0.01

Target Compounds Qvalue

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

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