

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009760.D
 Acq On : 11 Apr 2022 20:27
 Operator : CG/JU
 Sample : PB144015BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK015

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022

Quant Time: Apr 12 01:53:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	81196	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.781	136	359584	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.598	164	265875	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.351	188	593766	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.433	240	645358	20.000	ng/u1	-0.01
88) Perylene-d12	23.904	264	657835	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	13349m	7.337	ng/uL	0.00
4) Pyridine-d5	3.805	84	176483	35.183	ng/u1	0.00
7) Phenol-d5	7.122	99	229498	32.194	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	157442	37.099	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.499	132	185900	34.572	ng/u1	0.00
15) 4-Methylphenol-d8	8.669	113	203497	34.197	ng/u1	-0.01
21) Nitrobenzene-d5	9.128	128	101082	38.982	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.852	143	107319	37.750	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	194284	32.659	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.904	131	301966	35.639	ng/u1	-0.01
46) Dimethylphthalate-d6	14.010	166	811793	38.870	ng/u1	-0.01
49) Acenaphthylene-d8	14.292	160	841996	34.030	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.775	143	131618	34.750	ng/u1	-0.01
60) Fluorene-d10	15.592	176	664367	37.815	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	116715	33.181	ng/u1	-0.01
73) Anthracene-d10	17.451	188	1093783	38.474	ng/u1	-0.01
81) Pyrene-d10	19.680	212	1351208	38.336	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.745	264	1347397	39.331	ng/u1	-0.02

Target Compounds Qvalue

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

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