

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009759.D
 Acq On : 11 Apr 2022 19:53
 Operator : CG/JU
 Sample : PB144014BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK014

Manual Integrations
 APPROVED

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

Quant Time: Apr 12 01:53:02 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	113852	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.781	136	502206	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.598	164	368290	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.351	188	822888	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.433	240	843543	20.000	ng/u1	-0.01
88) Perylene-d12	23.904	264	772734	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	17951m	7.037	ng/uL	0.00
4) Pyridine-d5	3.805	84	243344	34.597	ng/u1	0.00
7) Phenol-d5	7.122	99	311775	31.192	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	211169	35.487	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.499	132	257018	34.088	ng/u1	0.00
15) 4-Methylphenol-d8	8.669	113	273244	32.747	ng/u1	-0.01
21) Nitrobenzene-d5	9.128	128	136034	37.562	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.852	143	150306	37.856	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	263240	31.683	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.905	131	410272	34.670	ng/u1	-0.01
46) Dimethylphthalate-d6	14.010	166	1074202	37.132	ng/u1	-0.01
49) Acenaphthylene-d8	14.293	160	1140135	33.266	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.781	143	181747	34.641	ng/u1	0.00
60) Fluorene-d10	15.593	176	877481	36.056	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.704	200	156695	32.144	ng/u1	0.00
73) Anthracene-d10	17.451	188	1459562	37.046	ng/u1	-0.01
81) Pyrene-d10	19.681	212	1744420	37.864	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.745	264	1547328	38.451	ng/u1	-0.02

Target Compounds Qvalue

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

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