

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010481.D
 Acq On : 23 May 2022 17:43
 Operator : CG/JU
 Sample : PB144963BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK963

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/24/2022
 Supervised By :Yogesh Patel 05/26/2022

Quant Time: May 23 23:20:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	159417	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.693	136	734494	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.522	164	499136	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.275	188	1061699	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.357	240	926470	20.000	ng/u1	# 0.00
88) Perylene-d12	23.774	264	696912	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	18738m	5.010	ng/uL	0.00
4) Pyridine-d5	3.758	84	283616	26.707	ng/u1	0.00
7) Phenol-d5	7.058	99	360124	26.915	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	251860	28.016	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.422	132	293966m	28.712	ng/u1	-0.01
15) 4-Methylphenol-d8	8.599	113	318004	27.966	ng/u1	-0.01
21) Nitrobenzene-d5	9.046	128	162735	28.628	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.769	143	174541	28.973	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	295931	25.980	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.828	131	438831m	25.910	ng/u1	-0.01
46) Dimethylphthalate-d6	13.934	166	1030548	28.023	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	1190745	28.071	ng/u1	0.00
54) 4-Nitrophenol-d4	14.722	143	133573	20.406	ng/u1	0.00
60) Fluorene-d10	15.516	176	887264	27.733	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	125925	20.280	ng/u1	0.00
73) Anthracene-d10	17.375	188	1390059	28.914	ng/u1	0.00
81) Pyrene-d10	19.604	212	1673722	34.017	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.621	264	1089374	31.130	ng/u1	-0.01

Target Compounds Qvalue

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

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