

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051022\
 Data File : BP010269.D
 Acq On : 10 May 2022 17:25
 Operator : CG/JU
 Sample : PB144591BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK591

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/11/2022
 Supervised By :mohammad ahmed 05/12/2022

Quant Time: May 11 01:17:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 10 15:59:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	115702	20.000	ng/u1	0.00
20) Naphthalene-d8	10.710	136	499152	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.539	164	359245	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.292	188	819693	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.374	240	847494	20.000	ng/u1	#-0.01
88) Perylene-d12	23.810	264	857497	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	18168m	6.671	ng/uL	0.00
4) Pyridine-d5	3.769	84	235015	31.009	ng/u1	0.00
7) Phenol-d5	7.075	99	289528	26.942	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	205637	30.933	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	237846m	30.053	ng/u1	0.00
15) 4-Methylphenol-d8	8.616	113	245752	27.816	ng/u1	0.00
21) Nitrobenzene-d5	9.069	128	123953	30.971	ng/u1	0.00
24) 2-Nitrophenol-d4	9.793	143	137192	30.064	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	229334	27.612	ng/u1	0.00
31) 4-Chloroaniline-d4	10.845	131	350379	28.575	ng/u1	-0.01
46) Dimethylphthalate-d6	13.951	166	937822	32.768	ng/u1	0.00
49) Acenaphthylene-d8	14.234	160	987618	28.944	ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	135502m	24.683	ng/u1	-0.01
60) Fluorene-d10	15.533	176	768775	31.822	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.651	200	147049	26.197	ng/u1	-0.01
73) Anthracene-d10	17.392	188	1266902	31.562	ng/u1	-0.01
81) Pyrene-d10	19.627	212	1567495	33.470	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.657	264	1516415	33.160	ng/u1	-0.01

Target Compounds Qvalue

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

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