

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010143.D
 Acq On : 29 Apr 2022 10:23
 Operator : CG/JU
 Sample : PB144447BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK447

Quant Time: Apr 29 10:55:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 05/02/2022
 Supervised By :Yogesh Patel 05/05/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	112440	20.000	ng/u1	#-0.01
20) Naphthalene-d8	10.734	136	441144	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.563	164	265802	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.316	188	506219	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.398	240	429543m	20.000	ng/u1	-0.01
88) Perylene-d12	23.845	264	418690	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	20331	7.885	ng/uL	0.00
4) Pyridine-d5	3.775	84	249312	34.370	ng/u1	-0.01
7) Phenol-d5	7.087	99	335327	32.787	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.252	67	234101	38.146	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.458	132	283346	37.409	ng/u1	-0.01
15) 4-Methylphenol-d8	8.634	113	278085	32.817	ng/u1	-0.02
21) Nitrobenzene-d5	9.087	128	137319	39.138	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.810	143	151089	39.429	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.352	165	256787	34.423	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.863	131	363563	33.990	ng/u1	-0.02
46) Dimethylphthalate-d6	13.969	166	814276	39.382	ng/u1	-0.01
49) Acenaphthylene-d8	14.251	160	944640	37.749	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.751	143	101293m	26.611	ng/u1	-0.01
60) Fluorene-d10	15.557	176	686698	38.609	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	107367m	33.543	ng/u1	0.00
73) Anthracene-d10	17.416	188	988724	40.372	ng/u1	0.00
81) Pyrene-d10	19.645	212	1089812	44.613	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.686	264	935339	42.612	ng/u1	-0.01

Target Compounds Qvalue

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

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