

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010158.D
 Acq On : 29 Apr 2022 18:57
 Operator : CG/JU
 Sample : PB144448BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK448

Manual Integrations
 APPROVED

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

Quant Time: Apr 30 00:32:09 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	144314	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.728	136	584216	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.557	164	362372	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.310	188	720416	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.398	240	638494	20.000	ng/u1	#-0.01
88) Perylene-d12	23.845	264	609510	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	20444	6.177	ng/uL	-0.01
4) Pyridine-d5	3.775	84	258745	27.792	ng/u1	-0.01
7) Phenol-d5	7.093	99	348213	26.527	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	241737	30.691	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.458	132	289710	29.801	ng/u1	-0.01
15) 4-Methylphenol-d8	8.634	113	289721	26.639	ng/u1	-0.02
21) Nitrobenzene-d5	9.087	128	144062	31.004	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.810	143	159194	31.370	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.351	165	262944	26.616	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.863	131	376817	26.602	ng/u1	-0.02
46) Dimethylphthalate-d6	13.969	166	862247	30.589	ng/u1	-0.01
49) Acenaphthylene-d8	14.251	160	1001946	29.369	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.757	143	115747m	22.304	ng/u1	0.00
60) Fluorene-d10	15.551	176	729068	30.067	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.669	200	113407	24.896	ng/u1	-0.01
73) Anthracene-d10	17.410	188	1102737	31.640	ng/u1	-0.01
81) Pyrene-d10	19.645	212	1248113	34.372	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1062870	33.263	ng/u1	-0.01

Target Compounds Qvalue

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

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