

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010837.D
 Acq On : 25 Jun 2022 20:41
 Operator : CG/JU
 Sample : PB145745BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK745

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/27/2022
 Supervised By :mohammad ahmed 07/01/2022

Quant Time: Jun 26 22:54:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	548739	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	2283350	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.375	164	1266521	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.133	188	2281383	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1833868	20.000	ng/u1	#-0.01
88) Perylene-d12	23.615	264	1955381	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	79371	5.472	ng/uL	0.00
4) Pyridine-d5	3.611	84	1036623	26.846	ng/u1	0.00
7) Phenol-d5	6.899	99	1245513	28.673	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	791153	27.849	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	1049775	29.129	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	1008944	29.382	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	493865	33.976	ng/u1	0.00
24) 2-Nitrophenol-d4	9.599	143	473915	39.408	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.140	165	846963	28.504	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	1407437	29.017	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	2656909	30.625	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	3146776	28.419	ng/u1	0.00
54) 4-Nitrophenol-d4	14.587	143	361714m	26.693	ng/u1	0.00
60) Fluorene-d10	15.369	176	2182321	29.312	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.492	200	173937	23.012	ng/u1	0.00
73) Anthracene-d10	17.233	188	3094874	30.491	ng/u1	-0.01
81) Pyrene-d10	19.486	212	3160948	31.638	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.463	264	3029765	31.355	ng/u1	-0.01

Target Compounds Qvalue

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

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