

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015870.D
 Acq On : 01 Jul 2023 06:24
 Operator : MA/JU
 Sample : PB153698BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK698

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/01/2023
 Supervised By :mohammad ahmed 07/01/2023

Quant Time: Jul 01 07:03:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 30 22:30:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.052	152	306013	20.000	ng/u1	0.00
20) Naphthalene-d8	10.875	136	1393861	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.698	164	875464	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.463	188	1859325	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.557	240	1353530	20.000	ng/u1	# 0.00
88) Perylene-d12	24.098	264	1401662	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	40315	4.740	ng/uL	0.00
4) Pyridine-d5	3.817	84	591263	27.577	ng/u1	0.00
7) Phenol-d5	7.228	99	732838	27.989	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.375	67	497277	30.698	ng/u1	0.00
11) 2-Chlorophenol-d4	7.581	132	588484	29.774	ng/u1	0.00
15) 4-Methylphenol-d8	8.781	113	616530	29.807	ng/u1	0.00
21) Nitrobenzene-d5	9.240	128	298741	28.045	ng/u1	0.00
24) 2-Nitrophenol-d4	9.969	143	323587	26.027	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.528	165	507086	22.888	ng/u1	0.01
31) 4-Chloroaniline-d4	11.040	131	817621	28.729	ng/u1	0.00
46) Dimethylphthalate-d6	14.110	166	1868803	27.618	ng/u1	0.00
49) Acenaphthylene-d8	14.398	160	2131316	28.019	ng/u1	0.00
54) 4-Nitrophenol-d4	14.981	143	224333m	25.122	ng/u1	0.03
60) Fluorene-d10	15.698	176	1574289	28.255	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	185142	16.191	ng/u1	0.00
73) Anthracene-d10	17.563	188	2494151	29.367	ng/u1	0.00
81) Pyrene-d10	19.792	212	2644839	29.925	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.927	264	2054402	29.055	ng/u1	0.00

Target Compounds Qvalue

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

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