

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
 Data File : BP020841.D
 Acq On : 08 Jul 2024 21:35
 Operator : MA/JU
 Sample : PB161754BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK754

Quant Time: Jul 08 23:06:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 03 15:01:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	36051	20.000	ng/u1	0.00
20) Naphthalene-d8	10.498	136	347262	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.357	164	416679	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.174	188	1010406	20.000	ng/u1	0.01
79) Chrysene-d12	21.621	240	989278	20.000	ng/u1	0.00
88) Perylene-d12	24.968	264	1113929	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.263	96	20982	24.992	ng/uL	0.00
4) Pyridine-d5	3.675	84	295222	120.657	ng/u1	0.00
7) Phenol-d5	6.916	99	375045	126.839	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	228456	121.874	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.263	132	307940	127.428	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	304913	126.969	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	152796	59.597	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.592	143	174934	68.182	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	293083	55.456	ng/u1	0.00
31) 4-Chloroaniline-d4	10.639	131	427551	57.727	ng/u1	0.00
46) Dimethylphthalate-d6	13.763	166	1010713	34.887	ng/u1	0.00
49) Acenaphthylene-d8	14.045	160	1078586	31.200	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	141096	31.875	ng/u1	0.03
60) Fluorene-d10	15.374	176	829777	34.036	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.510	200	143422	28.167	ng/u1	0.01
73) Anthracene-d10	17.274	188	1340195	29.554	ng/u1	0.01
81) Pyrene-d10	19.657	212	1633957	27.493	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.733	264	1622607	29.945	ng/u1	0.00

Target Compounds	Qvalue

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

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