

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041124\
 Data File : BM045140.D
 Acq On : 11 Apr 2024 18:44
 Operator : MA/JU
 Sample : PB160160BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK160

Quant Time: Apr 11 22:56:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 11 18:54:49 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	53406	20.000	ng/u1	0.00
20) Naphthalene-d8	10.392	136	229168	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.263	164	156534	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.027	188	338716	20.000	ng/u1	0.00
79) Chrysene-d12	21.233	240	285947	20.000	ng/u1	0.00
88) Perylene-d12	23.450	264	318504	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	7612	5.322	ng/uL	0.00
4) Pyridine-d5	3.605	84	111288	28.615	ng/u1	0.00
7) Phenol-d5	6.798	99	131058	28.947	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	88510	32.813	ng/u1	0.00
11) 2-Chlorophenol-d4	7.157	132	114578	32.302	ng/u1	0.00
15) 4-Methylphenol-d8	8.328	113	110465	31.171	ng/u1	0.00
21) Nitrobenzene-d5	8.781	128	57401	32.609	ng/u1	0.00
24) 2-Nitrophenol-d4	9.492	143	63672	34.269	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	103441	30.365	ng/u1	0.00
31) 4-Chloroaniline-d4	10.551	131	160686	29.627	ng/u1	0.00
46) Dimethylphthalate-d6	13.692	166	412250	37.824	ng/u1	0.00
49) Acenaphthylene-d8	13.957	160	430084	33.402	ng/u1	0.00
54) 4-Nitrophenol-d4	14.510	143	54370	23.903	ng/u1	0.01
60) Fluorene-d10	15.268	176	333994	34.373	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	47996	23.899	ng/u1	0.00
73) Anthracene-d10	17.127	188	531639	34.891	ng/u1	0.00
81) Pyrene-d10	19.433	212	613705	43.587	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.309	264	567272	36.510	ng/u1	0.00

Target Compounds Qvalue

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

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