

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
 Data File : BM045023.D
 Acq On : 08 Apr 2024 01:07
 Operator : MA/JU
 Sample : PB160066BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK066

Quant Time: Apr 08 04:20:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 01:45:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	84104	20.000	ng/u1	0.00
20) Naphthalene-d8	10.398	136	367430	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.274	164	233454	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.033	188	484610	20.000	ng/u1	-0.01
79) Chrysene-d12	21.239	240	429112	20.000	ng/u1	-0.01
88) Perylene-d12	23.462	264	516564	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	13798	6.126	ng/uL	0.00
4) Pyridine-d5	3.605	84	203272	33.189	ng/u1	0.00
7) Phenol-d5	6.810	99	243078	34.093	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	147839	34.803	ng/u1	0.00
11) 2-Chlorophenol-d4	7.163	132	205324	36.757	ng/u1	0.00
15) 4-Methylphenol-d8	8.334	113	198797	35.621	ng/u1	0.00
21) Nitrobenzene-d5	8.787	128	99286	35.179	ng/u1	0.00
24) 2-Nitrophenol-d4	9.498	143	110811	37.198	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.034	165	187611	34.350	ng/u1	0.00
31) 4-Chloroaniline-d4	10.557	131	277377	31.898	ng/u1	0.00
46) Dimethylphthalate-d6	13.704	166	564408	34.723	ng/u1	0.00
49) Acenaphthylene-d8	13.969	160	680027	35.413	ng/u1	0.00
54) 4-Nitrophenol-d4	14.516	143	90414	26.652	ng/u1	0.00
60) Fluorene-d10	15.274	176	516388	35.634	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.421	200	66466	23.132	ng/u1	0.00
73) Anthracene-d10	17.133	188	807848	37.057	ng/u1	-0.01
81) Pyrene-d10	19.439	212	914496	43.281	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.321	264	940097	37.306	ng/u1	-0.01

Target Compounds Qvalue

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

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