

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060721\
 Data File : BM030308.D
 Acq On : 07 Jun 2021 14:39
 Operator : CG/JU
 Sample : PB136692BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK592

Quant Time: Jun 07 15:10:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 07 09:55:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	189213	20.000	ng/u1	0.00
20) Naphthalene-d8	10.645	136	769394	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.474	164	474231	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.210	188	941129	20.000	ng/u1	0.00
79) Chrysene-d12	21.368	240	851898	20.000	ng/u1	0.00
88) Perylene-d12	23.656	264	873073	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.316	96	32059	6.470	ng/uL	0.00
4) Pyridine-d5	3.728	84	354668	27.239	ng/u1	0.00
7) Phenol-d5	7.004	99	466479	28.366	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	309540	29.164	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	384400	32.321	ng/u1	0.00
15) 4-Methylphenol-d8	8.545	113	361384	27.515	ng/u1	0.00
21) Nitrobenzene-d5	9.004	128	173098	32.867	ng/u1	0.00
24) 2-Nitrophenol-d4	9.728	143	169434	32.392	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	352602	30.947	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	500106	28.712	ng/u1	0.00
46) Dimethylphthalate-d6	13.880	166	1085462	33.353	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	1343690	33.049	ng/u1	0.00
54) 4-Nitrophenol-d4	14.651	143	157843	25.911	ng/u1	0.00
60) Fluorene-d10	15.463	176	968496	33.241	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	132406	24.373	ng/u1	0.00
73) Anthracene-d10	17.310	188	1456828	34.034	ng/u1	0.00
81) Pyrene-d10	19.592	212	1567392	35.152	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.509	264	1554942	35.717	ng/u1	0.00

Target Compounds Qvalue

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

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