

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
 Data File : BM030839.D
 Acq On : 07 Jul 2021 03:38
 Operator : CG/JU
 Sample : PB137437BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK437

Quant Time: Jul 07 05:02:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 07 04:59:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	121450	20.000	ng/u1	0.00
20) Naphthalene-d8	10.545	136	507913	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.392	164	330802	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.133	188	672802	20.000	ng/u1	0.00
79) Chrysene-d12	21.303	240	515038	20.000	ng/u1	0.00
88) Perylene-d12	23.544	264	491671	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	20800	6.692	ng/uL	0.00
4) Pyridine-d5	3.669	84	273802	33.687	ng/u1	0.00
7) Phenol-d5	6.934	99	352679	33.234	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.098	67	233129	34.105	ng/u1	0.00
11) 2-Chlorophenol-d4	7.298	132	292373	35.910	ng/u1	0.00
15) 4-Methylphenol-d8	8.463	113	285813	33.607	ng/u1	0.00
21) Nitrobenzene-d5	8.922	128	137326	36.352	ng/u1	0.00
24) 2-Nitrophenol-d4	9.639	143	154639	37.412	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	274816	34.549	ng/u1	0.00
31) 4-Chloroaniline-d4	10.686	131	393733	34.856	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	908247	39.218	ng/u1	0.00
49) Acenaphthylene-d8	14.086	160	1110879	37.974	ng/u1	0.00
54) 4-Nitrophenol-d4	14.592	143	130520	32.384	ng/u1	0.00
60) Fluorene-d10	15.386	176	781378	38.316	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	118581	27.487	ng/u1	0.00
73) Anthracene-d10	17.233	188	1205685	38.456	ng/u1	0.00
81) Pyrene-d10	19.521	212	1214175	42.745	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.403	264	972182	38.306	ng/u1	0.00

Target Compounds	Qvalue

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

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