

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
 Data File : BM032442.D
 Acq On : 12 Oct 2021 05:03
 Operator : CG/JU
 Sample : PB139788BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK788

Quant Time: Oct 12 05:43:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 11 11:06:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.613	152	222188	20.000	ng/u1	0.00
20) Naphthalene-d8	10.401	136	1030865	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.260	164	724440	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.989	188	1521268	20.000	ng/u1	0.00
79) Chrysene-d12	21.148	240	1468498	20.000	ng/u1	-0.01
88) Perylene-d12	23.294	264	1504189	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.955	96	36791	6.639	ng/uL	0.00
4) Pyridine-d5	3.384	84	520394	33.241	ng/u1	0.00
7) Phenol-d5	6.801	99	638733	33.807	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.937	67	381098	35.130	ng/u1	0.00
11) 2-Chlorophenol-d4	7.148	132	513695	35.610	ng/u1	0.00
15) 4-Methylphenol-d8	8.337	113	540650	35.429	ng/u1	-0.01
21) Nitrobenzene-d5	8.766	128	266559	36.649	ng/u1	0.00
24) 2-Nitrophenol-d4	9.489	143	302133	38.537	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.031	165	525369	33.662	ng/u1	0.00
31) 4-Chloroaniline-d4	10.536	131	793967	34.490	ng/u1	0.00
46) Dimethylphthalate-d6	13.666	166	1845158	36.803	ng/u1	0.00
49) Acenaphthylene-d8	13.948	160	2204992	36.203	ng/u1	0.00
54) 4-Nitrophenol-d4	14.466	143	318544	32.353	ng/u1	-0.01
60) Fluorene-d10	15.248	176	1554463	36.559	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.366	200	273154	31.506	ng/u1	0.00
73) Anthracene-d10	17.083	188	2443097	36.894	ng/u1	0.00
81) Pyrene-d10	19.371	212	2745134	37.158	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.159	264	2655254	35.250	ng/u1	-0.01

Target Compounds	Qvalue

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

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