

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121022\
 Data File : BM037945.D
 Acq On : 10 Dec 2022 22:08
 Operator : CG/JU
 Sample : PB149486BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK486

Quant Time: Dec 11 21:30:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 09 21:34:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.075	152	198070	20.000	ng/u1	0.00
20) Naphthalene-d8	10.898	136	759387	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.704	164	399607	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	729487	20.000	ng/u1	0.00
79) Chrysene-d12	21.615	240	729716	20.000	ng/u1	0.00
88) Perylene-d12	24.097	264	785146	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	23601	5.410	ng/uL	0.00
4) Pyridine-d5	3.857	84	298812	23.091	ng/u1	0.00
7) Phenol-d5	7.234	99	401083	24.919	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.398	67	246520	25.153	ng/u1	0.00
11) 2-Chlorophenol-d4	7.604	132	351355	26.611	ng/u1	0.00
15) 4-Methylphenol-d8	8.786	113	307484	24.474	ng/u1	0.00
21) Nitrobenzene-d5	9.251	128	164930	27.431	ng/u1	0.00
24) 2-Nitrophenol-d4	9.975	143	170221	27.016	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.516	165	299364	25.014	ng/u1	0.00
31) 4-Chloroaniline-d4	11.033	131	386786	23.413	ng/u1	0.00
46) Dimethylphthalate-d6	14.110	166	741696	26.034	ng/u1	-0.01
49) Acenaphthylene-d8	14.404	160	909973	26.134	ng/u1	0.00
54) 4-Nitrophenol-d4	14.898	143	93942	19.348	ng/u1	0.00
60) Fluorene-d10	15.692	176	654412	25.783	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	60430	15.494	ng/u1	0.00
73) Anthracene-d10	17.545	188	971884	28.546	ng/u1	0.00
81) Pyrene-d10	19.827	212	1126839	25.707	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.933	264	1144037	29.172	ng/u1	-0.01

Target Compounds Qvalue

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

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