

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121022\
 Data File : BM037948.D
 Acq On : 10 Dec 2022 23:57
 Operator : CG/JU
 Sample : N5927-13
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ89

Quant Time: Dec 11 21:30:36 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	164605	20.000	ng/u1	0.00
20) Naphthalene-d8	10.898	136	641421	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.704	164	348307	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	624684	20.000	ng/u1	0.00
79) Chrysene-d12	21.615	240	554584	20.000	ng/u1	0.00
88) Perylene-d12	24.097	264	615737	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	18397	5.075	ng/uL	0.00
4) Pyridine-d5	3.863	84	85071	7.910	ng/u1	0.00
7) Phenol-d5	7.228	99	92316	6.902	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.398	67	285091	35.003	ng/u1	0.00
11) 2-Chlorophenol-d4	7.604	132	304409	27.742	ng/u1	0.00
15) 4-Methylphenol-d8	8.786	113	167999	16.090	ng/u1	0.00
21) Nitrobenzene-d5	9.251	128	195644	38.524	ng/u1	0.00
24) 2-Nitrophenol-d4	9.975	143	191258	35.937	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.516	165	315992	31.259	ng/u1	0.00
31) 4-Chloroaniline-d4	11.033	131	391424	28.051	ng/u1	0.00
46) Dimethylphthalate-d6	14.116	166	991646	39.934	ng/u1	0.00
49) Acenaphthylene-d8	14.404	160	1158614	38.176	ng/u1	0.00
54) 4-Nitrophenol-d4	14.904	143	16472	3.892	ng/u1	0.00
60) Fluorene-d10	15.692	176	863185	39.017	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	80721	24.169	ng/u1	0.00
73) Anthracene-d10	17.545	188	1292649	44.338	ng/u1	0.00
81) Pyrene-d10	19.827	212	1437166	43.140	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.933	264	1400050	45.523	ng/u1	-0.01
Target Compounds						
2) 1,4-Dioxane	3.457	88	4474	1.176	ng/uL	91

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

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