

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040724\
 Data File : BM045020.D
 Acq On : 07 Apr 2024 22:43
 Operator : MA/JU
 Sample : P1973-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YCST6

Quant Time: Apr 08 04:14:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 08 01:45:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	78599	20.000	ng/u1	0.00
20) Naphthalene-d8	10.398	136	333446	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.274	164	207266	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.033	188	439174	20.000	ng/u1	-0.01
79) Chrysene-d12	21.244	240	403510	20.000	ng/u1	0.00
88) Perylene-d12	23.462	264	482505	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.204	96	8407	3.994	ng/uL	0.00
4) Pyridine-d5	3.616	84	51601	9.015	ng/u1	0.00
7) Phenol-d5	6.816	99	39010	5.854	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	137043	34.521	ng/u1	0.00
11) 2-Chlorophenol-d4	7.163	132	138144	26.463	ng/u1	0.00
15) 4-Methylphenol-d8	8.339	113	79929	15.325	ng/u1	0.00
21) Nitrobenzene-d5	8.786	128	94265	36.804	ng/u1	0.00
24) 2-Nitrophenol-d4	9.504	143	93370	34.538	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.033	165	159507	32.181	ng/u1	0.00
31) 4-Chloroaniline-d4	10.557	131	209202	26.510	ng/u1	0.00
46) Dimethylphthalate-d6	13.704	166	562308	38.964	ng/u1	0.00
49) Acenaphthylene-d8	13.968	160	655178	38.429	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	10391	3.450	ng/u1	0.03
60) Fluorene-d10	15.274	176	516103	40.114	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.421	200	74005	28.420	ng/u1	0.00
73) Anthracene-d10	17.133	188	883225	44.706	ng/u1	-0.01
81) Pyrene-d10	19.439	212	1053056	53.001	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.321	264	1098321	46.662	ng/u1	-0.01

Target Compounds	Qvalue

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

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