

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
 Data File : BM042902.D
 Acq On : 19 Nov 2023 12:36
 Operator : MA/JU
 Sample : PB157267BL
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 19 21:48:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 18 22:06:09 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	28175	20.000	ng/u1	0.00
20) Naphthalene-d8	10.622	136	110734	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.468	164	70523	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	157160	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.409	240	163945	20.000	ng/u1	0.00
88) Perylene-d12	23.762	264	214677	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	4959	5.523	ng/uL	0.00
4) Pyridine-d5	3.604	84	66236	29.655	ng/u1	0.00
7) Phenol-d5	7.104	99	204	0.074	ng/u1	0.13
9) Bis-(2-Chloroethyl)eth...	7.163	67	57890	30.044	ng/u1	0.00
11) 2-Chlorophenol-d4	7.334	132	56689	26.792	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	52614	24.369	ng/u1	0.00
21) Nitrobenzene-d5	9.157	128	142	0.146	ng/u1	0.14
24) 2-Nitrophenol-d4	9.728	143	27363	24.093	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	48467	22.951	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	55509	20.702	ng/u1	0.00
46) Dimethylphthalate-d6	13.892	166	172621	25.863	ng/u1	0.00
49) Acenaphthylene-d8	14.168	160	187115	26.201	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	4721	6.261	ng/u1	0.02
60) Fluorene-d10	15.462	176	150419	27.215	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.615	200	18396	15.453	ng/u1	0.00
73) Anthracene-d10	17.321	188	234345	26.607	ng/u1	0.00
81) Pyrene-d10	19.621	212	317606	29.584	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.609	264	353268	27.511	ng/u1	0.00

Target Compounds Qvalue

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

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