

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101123\
 Data File : BP017647.D
 Acq On : 11 Oct 2023 15:13
 Operator : MA/JU
 Sample : PB156228TB 10X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB228

Quant Time: Oct 12 02:22:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 10 04:45:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	98266	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	385424	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.622	164	244179	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	585533	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	594221	20.000	ng/u1	0.00
88) Perylene-d12	24.133	264	666191	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	17224	7.111	ng/uL	0.00
4) Pyridine-d5	3.716	84	214033	31.391	ng/u1	0.00
7) Phenol-d5	7.081	99	268504	32.311	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	171638	33.303	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	225609	34.514	ng/u1	0.00
15) 4-Methylphenol-d8	8.646	113	212144	31.693	ng/u1	0.00
21) Nitrobenzene-d5	9.134	128	103866	34.973	ng/u1	0.00
24) 2-Nitrophenol-d4	9.851	143	110948	34.014	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	200274	31.614	ng/u1	0.00
31) 4-Chloroaniline-d4	10.934	131	289568	32.093	ng/u1	0.00
46) Dimethylphthalate-d6	14.039	166	704584	36.343	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	727310	34.457	ng/u1	0.00
54) 4-Nitrophenol-d4	14.875	143	75883	25.825	ng/u1	0.01
60) Fluorene-d10	15.628	176	611387	36.060	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	85192	24.529	ng/u1	0.00
73) Anthracene-d10	17.516	188	1035524	36.851	ng/u1	0.00
81) Pyrene-d10	19.769	212	1299224	38.331	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.962	264	1283028	37.148	ng/u1	0.00

Target Compounds Qvalue

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

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