

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
 Data File : BP016128.D
 Acq On : 09 Jul 2023 12:16
 Operator : MA/JU
 Sample : 03332-20
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 09 22:40:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 08 10:11:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	297331	20.000	ng/ul	0.00
20) Naphthalene-d8	10.851	136	1294847	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.675	164	824741	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	1725307	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.551	240	1480344	20.000	ng/ul	# 0.01
88) Perylene-d12	24.080	264	1372391	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	23806	2.881	ng/uL	0.00
4) Pyridine-d5	3.805	84	244927	11.757	ng/ul	0.00
7) Phenol-d5	7.240	99	319145	12.545	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.357	67	191170	12.146	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	291990	15.205	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	290370	14.448	ng/ul	0.02
21) Nitrobenzene-d5	9.240	128	129592	13.096	ng/ul	0.02
24) 2-Nitrophenol-d4	9.987	143	149287	12.926	ng/ul	0.04
28) 2,4-Dichlorophenol-d3	10.587	165	187277	9.099	ng/ul	0.08
31) 4-Chloroaniline-d4	11.069	131	327600	12.391	ng/ul	0.05
46) Dimethylphthalate-d6	14.104	166	1068246	16.758	ng/ul	0.02
49) Acenaphthylene-d8	14.369	160	1152839	16.088	ng/ul	0.00
54) 4-Nitrophenol-d4	14.957	143	19	0.002	ng/ul	0.00
60) Fluorene-d10	15.686	176	890316	16.962	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.916	200	44777	4.220	ng/ul	0.06
73) Anthracene-d10	17.551	188	1339223	16.993	ng/ul	0.00
81) Pyrene-d10	19.780	212	1573352	16.277	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.909	264	1140407	16.473	ng/ul	0.00

Target Compounds	Qvalue

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

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