

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009890.D
 Acq On : 16 Apr 2022 08:26
 Operator : CG/JU
 Sample : N2421-16
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J76

Quant Time: Apr 18 01:10:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	129924	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	600030	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	440891	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	1012123	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	1034617	20.000	ng/ul	# 0.00
88) Perylene-d12	23.886	264	962407	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	12476	4.286	ng/uL	0.00
4) Pyridine-d5	3.799	84	81580	10.164	ng/ul	0.00
7) Phenol-d5	7.111	99	72503	6.356	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	218683	32.203	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	201786	23.452	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	148923	15.640	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	141796	32.770	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	148321	31.266	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	256635	25.853	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	361976	25.602	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	1104538	31.893	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1203975	29.344	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	38025	6.054	ng/ul	0.00
60) Fluorene-d10	15.581	176	961592	33.006	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	172786	28.818	ng/ul	0.00
73) Anthracene-d10	17.439	188	1717818	35.448	ng/ul	0.00
81) Pyrene-d10	19.669	212	2129121	37.680	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1921537	38.340	ng/ul	0.00

Target Compounds Qvalue

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

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