

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009908.D
 Acq On : 18 Apr 2022 11:06
 Operator : CG/JU
 Sample : PB144126BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK126

Quant Time: Apr 19 00:53:11 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	7.952	152	118203	20.000	ng/u1	0.00
20) Naphthalene-d8	10.763	136	517173	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	379694	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.334	188	875962	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.416	240	869185	20.000	ng/u1	# 0.00
88) Perylene-d12	23.874	264	777082	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	16163	6.103	ng/uL	0.00
4) Pyridine-d5	3.793	84	219792	30.099	ng/u1	0.00
7) Phenol-d5	7.111	99	278369	26.824	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	186637	30.210	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	227774	29.098	ng/u1	0.00
15) 4-Methylphenol-d8	8.658	113	241767	27.908	ng/u1	0.00
21) Nitrobenzene-d5	9.110	128	121718	32.637	ng/u1	0.00
24) 2-Nitrophenol-d4	9.840	143	130821	31.995	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	236016	27.585	ng/u1	0.00
31) 4-Chloroaniline-d4	10.887	131	361196	29.639	ng/u1	0.00
46) Dimethylphthalate-d6	13.993	166	900999	30.209	ng/u1	0.00
49) Acenaphthylene-d8	14.275	160	979563	27.722	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.763	143	129589	23.958	ng/u1	0.00
60) Fluorene-d10	15.575	176	768798	30.641	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	121114	23.340	ng/u1	0.00
73) Anthracene-d10	17.434	188	1268972	30.257	ng/u1	0.00
81) Pyrene-d10	19.663	212	1567576	33.022	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.716	264	1277992	31.581	ng/u1	-0.01

Target Compounds Qvalue

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

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