

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009939.D
 Acq On : 19 Apr 2022 05:03
 Operator : CG/JU
 Sample : PB144126BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK126

Quant Time: Apr 19 06:34:32 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	7.951	152	167737	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	738128	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.580	164	533634	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.333	188	1190902	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.415	240	1151148	20.000	ng/u1	# 0.00
88) Perylene-d12	23.880	264	987533	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	23433	6.235	ng/uL	0.00
4) Pyridine-d5	3.793	84	312324	30.140	ng/u1	0.00
7) Phenol-d5	7.104	99	405795	27.556	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	270720	30.879	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	328275	29.552	ng/u1	0.00
15) 4-Methylphenol-d8	8.651	113	350557	28.516	ng/u1	-0.01
21) Nitrobenzene-d5	9.110	128	176456	33.151	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	194417	33.315	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	336083	27.522	ng/u1	0.00
31) 4-Chloroaniline-d4	10.886	131	509617	29.300	ng/u1	0.00
46) Dimethylphthalate-d6	13.986	166	1261919	30.105	ng/u1	-0.01
49) Acenaphthylene-d8	14.275	160	1404596	28.284	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.763	143	192495	25.322	ng/u1	0.00
60) Fluorene-d10	15.574	176	1063726	30.166	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	183227	25.971	ng/u1	0.00
73) Anthracene-d10	17.433	188	1713801	30.056	ng/u1	0.00
81) Pyrene-d10	19.662	212	2076560	33.029	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.721	264	1642902	31.946	ng/u1	0.00

Target Compounds Qvalue

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

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