

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006925.D
 Acq On : 10 Sep 2021 06:03
 Operator : CG/JU
 Sample : M3651-14
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBKM8

Quant Time: Sep 10 06:43:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 09 11:07:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	286286	20.000	ng/u1	0.00
20) Naphthalene-d8	10.734	136	1171234	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.557	164	723456	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.304	188	1510135	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.386	240	1554826	20.000	ng/u1	-0.01
88) Perylene-d12	23.821	264	1629602	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	19513	2.712	ng/uL	0.00
4) Pyridine-d5	3.793	84	164282	8.763	ng/u1	0.01
7) Phenol-d5	7.093	99	171198	8.136	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	450058	37.243	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	472478	27.815	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	303274	18.354	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	280974	39.606	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	251668	36.033	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	496874	30.171	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	726460	31.898	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	1832558	39.110	ng/u1	0.00
49) Acenaphthylene-d8	14.251	160	2228057	37.545	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	41360	5.081	ng/u1	0.00
60) Fluorene-d10	15.551	176	1662967	40.213	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	197001	29.489	ng/u1	0.00
73) Anthracene-d10	17.404	188	2764211	44.213	ng/u1	-0.01
81) Pyrene-d10	19.639	212	3201751	43.970	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.662	264	3289098	41.837	ng/u1	-0.02

Target Compounds Qvalue

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

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