

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG042420\  
 Data File : BG045173.D  
 Acq On : 24 Apr 2020 23:03  
 Operator : CG/JU  
 Sample : L2118-01 5X  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

**Instrument :**  
 BNA\_G  
**ClientSampleId :**  
 EO-02-042320

**Manual Integrations**  
**APPROVED**  
 mohammad  
 4/27/2020 3:05:52 PM

Quant Time: Apr 25 01:18:30 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Apr 23 13:58:55 2020  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	72205	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	281438	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	206013	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	482832	20.00	ng	0.00
76) Chrysene-d12	21.64	240	461931	20.00	ng	-0.01
87) Perylene-d12	24.82	264	492905	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	53516	12.24	ng	0.00
7) Phenol-d6	7.16	99	80382	12.37	ng	0.00
23) Nitrobenzene-d5	9.15	82	55631	9.02	ng	-0.01
42) 2,4,6-Tribromophenol	16.13	330	41149	12.93	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	134478	9.57	ng	0.00
79) Terphenyl-d14	19.99	244	245416	11.58	ng	0.00

Target Compounds

						Qvalue
71) Phenanthrene	17.42	178	67296	2.712	ng	96
75) Fluoranthene	19.43	202	375116	7.756	ng	96
78) Pyrene	19.79	202	414447	13.014	ng	95
81) Benzo(a)anthracene	21.62	228	243136	8.121	ng	99
83) Chrysene	21.69	228	230383	8.009	ng	97
86) Indeno(1,2,3-cd)pyrene	28.41	276	180309	4.721	ng	# 56
88) Benzo(b)fluoranthene	23.82	252	358533	11.612	ng	94
89) Benzo(k)fluoranthene	23.87	252	113592m	3.869	ng	
90) Benzo(a)pyrene	24.66	252	214888	7.372	ng	96
92) Benzo(a,h,i)perylene	29.54	276	175581	5.962	ng	98

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

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